Amendments to the Claims

1. (Currently Amended). A compound having a formula I,

$$Z \xrightarrow{A_3} Y \xrightarrow{R^2} A_2 \xrightarrow{E_2} E_1 \xrightarrow{R^3} E_4 \xrightarrow{R_4} R_5$$

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A₁ is: a bond, CH₂. O or S, and wherein A₂ and R⁴ or A₂ and R⁵ together being a 3- to 6-

membered carbocyclyl when A₁ is a carbon;

A2 and A3 are independently: CH2, O or S;

E₁, E₂, E₃, E₄ and E₅ are each CH or substituted carbon bearing A₂ and R³; or at least one of E₁, E₂, E₃, E₄ and E₅ is nitrogen and each of others being CH or substituted carbon bearing A₂ and R³;

Q is: -C(O)OR⁶, or R^{6A};

Y is: a bond, or C1-C6 alkyl-or C3-C6 eycloalkyl;

- Z is: a) phenyl-T-pyridinyl or naphthyl-T-pyridinyl each optionally substituted with one or more R⁷; wherein T is a single bond, C or O; aryl:
 - a 5- to 10-membered heteroaryl wherein the heteroaryl containing at least one heteroatom selected from N, O or S,
 - bi aryl, wherein biaryl being defined as aryl substituted with another aryl or aryl substituted with heteroaryl, or
 - bi-heteroaryl, wherein bi-heteroaryl being defined as heteroaryl substituted with another heteroaryl, or heteroaryl substituted with aryl, and wherein aryl, heteroaryl, bi-aryl and bi-heteroaryl being optionally substituted with one or more groups independently selected from R²;

n is: 1, 2, 3, 4, 5 or 6

p is: 1 or 2:

r is: 1, 2, 3, or 4:

R1 and R2 are each independently:

hydrogen,

haloalkyl,

C1-C6 alkyl,

(CH₂)_nC₃-C₈-cycloalkyl, or

R* and R2 form a 4 to 8 membered nonaromatic carbocyclic ring; and wherein at least one of R* and R2 is alkyl or cycloalkyl, and;

R3 is: hydrogen,

nitro.

cvano.

hydroxyl,

halo.

haloalkyl,

haloalkyloxy,

aryloxy,

C1-C6 alkyl,

C1-C6 alkoxy, or

C3-C8 cycloalkyl;

R4 and R5 are each independently: hydrogen or C4 C6 alkyl;

R6 is: hydrogen, C1-C6 alkyl or aminoalkyl;

R^{6A}-is:-carboxamide, sulfonamide, acylsulfonamide, tetrazole,

R7 is: hydrogen,

oxo,

nitro.

cyano,

hydroxyl,

halo,

haloalkyl,

haloalkyloxy,

aryloxy,

arylalkyl,

aminoalkyl,

C1-C6 alkyl,

C1-C6 alkoxy,

(CH₂)_nC₃-C₈ cycloalkyl,

 $C(O)R^9$

 $C(O)OR^9$,

 $C(=NOR^8)R^9$,

CR8(OH)R9,

 $C[=C(R^8)_2]R^9$,

OR9,

SR⁹ or

 $S(O)_pR^9$;

R8 is: hydrogen or C1-C6 alkyl; and

R9 is: hydrogen,

C1-C6 alkyl,

C3-C8 cycloalkyl,

aryl,

heteroaryl or

heterocyclyl,

wherein alkyl, cycloalkyl, aryl, heteroaryl or heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of:

hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1 - C_6 alkyl, C_1 - C_6 alkoxy and C_3 - C_8 cycloalkyl.

 (Currently Amended). The compound of Claim 1, wherein the compound is represented by a compound of formula II,

$$Z \xrightarrow{Q} A_2 \xrightarrow{(R^2)_r} A_1 \xrightarrow{Q} A_2$$

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A₁ is: a-bond, CH₂, O or S, and wherein A₄ and R⁴ or A₄ and R⁵ together being a 3 to 6-membered carbocyclyl when A₄ is a carbon;

A2 is: O or S or CH2;

Q is: $-C(O)OR^6$; or R^{6A} ;

Y is: a bond or; C1-C6 alkyl or C2-C6 eyeloalkyl;

Z is: phenyl-T-pyridinyl or naphthyl-T-pyridinyl each optionally substituted with one or more R⁷; wherein T is a single bond, C or O;a)—aryl;

 a 5 to 10 membered heteroaryl wherein the heteroaryl containing at least one heteroatom-selected from N, O or S;

```
e) bi-aryl, wherein biaryl being defined as aryl substituted with another aryl or aryl
               substituted with heteroaryl, or
            bi heterogryl, wherein bi heterogryl being defined as heterogryl substituted with
               another heteroaryl, or heteroaryl substituted with aryl, and
               wherein arvl, beteroarvl, bi arvl and bi beteroarvl being ontionally substituted with
               one or more groups independently selected from R2:
n is: 1, 2, 3, 4, 5 or 6
p is: 1 or 2:
r is: 1, 2, 3, or 4;
R1 and R2 are each independently:
       hydrogen,
       haloalkyl.
       C1-C6 alkylz
       (CH2),C2-C2-eveloalkyl, or
       R<sup>+</sup> and R<sup>2</sup> form a 4- to 8-membered nonaromatic carbocyclic ring; and
       wherein at least one of R<sup>1</sup> and R<sup>2</sup> is alkyl or eveloalkyl, and:
```

R³ is: hydrogen,
nitro,
cyano,
hydroxyl,
halo,
haloalkyl,
haloalkyloxy,
aryloxy,
C1-C6 alkyl,
C1-C6 alkoxy or
C1-C5 evcloalkyl;

R4 and R5 are each independently: hydrogen or C1-C6 alleyl;

R6 is: hydrogen, C1-C6 alkyl or aminoalkyl;

R^{6A} is: carboxamide, sulfonamide, acylsulfonamide, tetrazole,

R7 is: hydrogen,

oxo,

nitro.

cyano,

hydroxyl,

halo,

haloalkyl,

haloalkyloxy,

aryloxy,

arylalkyl,

aminoalkyl,

C₁-C₆ alkyl,

.

C1-C6 alkoxy,

(CH₂)_nC₃-C₈ cycloalkyl,

 $C(O)R^9$,

C(O)OR9,

 $C(=NOR^8)R^9$,

CR8(OH)R9,

 $C[=C(R^8)_2]R^9$,

OR9.

SR9 or

 $S(O)_pR^9$;

R8 is: hydrogen or C1-C6 alkyl; and

R9 is: hydrogen,

C1-C6 alkyl,

C3-C8 cycloalkyl,

arvl.

heteroaryl or

heterocyclyl,

wherein alkyl, cycloalkyl, aryl, heteroaryl or heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of:

hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1 - C_6 alkyl, C_1 - C_6 alkoxy and C_3 - C_8 cycloalkyl.

3. (<u>Currently Amended</u>). The compound of Claim 2, wherein Z is <u>an optionally</u> substituted phenyl or naphthyl, furanyl, imidazolyl, indolyl, oxazolyl, isoxazolyl, pyridyl, pyrrolyl, thiazolyl, thiazolyl, thiaphenyl, benzefuranyl, benzethiophenyl, benzeisoxazolyl, quinolinyl, isoquinolinyl or a structural formula selected from following:

wherein T is:

$$\begin{split} &a \ \text{bond,} \ \frac{(CH_2)_qO_{\tau} - O(CH_2)_{q^{\tau}} - C(O)(CH_2)_{q^{\tau}} - (CH_2)_qC(O)_{\tau} - (CH_2)_qS_{\tau} - S(CH_2)_{q^{\tau}} - S(O)_{p_{\tau}} \\ &- (C_1 - C_2 \ \text{all}_{N^2})_{1} - (CH_2)_qC(-CH_2)_{1} - C(-CH_2)_{1} - (CH_2)_qC(-NOH)_{\tau} \\ &- C(-NOH)(CH_2)_{q^{\tau}} - (CH_2)_qC(-NOCH_2)_{1} - C(-NOCH_2)_{1} - CH(OH)(CH_2)_{q^{\tau}} - OF_{\tau} \\ &- (CH_2)_qC(-NOCH_2)_{1} - (CH_2)_qC(-NOCH_2)_{2} - ($$

a is: 0, 1, 2 or 3; and

rings b to 1 rings g and h are each optionally substituted with one or more groups independently selected from the group consisting of:

hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl. S(O)₂R², C₁-C₆ alkyl. C₁-C₆ alkyl. C₁-C₆ alkyl. C₁-C₆ alkyl. C₁-C₆ alkyl.

4. (Canceled)

 (Withdrawn). The compound of Claim 2, wherein the compound is represented by structural formula IV.

$$R^{1}$$
 $COOR^{0}$
 $COOR^{0}$
 $COOR^{0}$
 $COOR^{0}$
 $COOR^{0}$

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein: A_1 and A_2 are respectively:

O and O.

CH2 and O,

CH2 and S.

O and S or

O 44114 15

S and O;

m is: 1 or 2;

R1 is: C1-C3 alkvl:

R3 is: hydrogen, halo or C1-C6 alkyl;

R6 and R9 are each independently; hydrogen or C1-C6 alkyl;

T is: a bond, -O-, -C(O)-, -S(O) -S(O)2-, -C(=CH₂)-, -C(=NOH)- or -CH(OH)-; and rings b and c are each optionally substituted with one or more groups independently selected from:

 $hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, \\ aminoalkyl, S(O)_2R^9, C_1-C_6 alkyl, C_1-C_6 alkoxy and (CH_2)_nC_3-C_8 cycloalkyl.$

 (Withdrawn). The compound of Claim 5, wherein the compound is represented by structural formula V,

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

T is: a bond, -O- or -C(O)-;

R1 is: methyl, ethyl or cyclopropyl;

R3 is: methyl or ethyl; and

rings b and c are each optionally substituted with one or more substituent independently selected from the group consisting of: hydrogen, Cl, Br, CF₃, OCF₃, methyl, ethyl, isopropyl, N(CH₃)₂, S(O)₂CH₃, methoxy and cyclopropyl.

 (Withdrawn). The compound of Claim 6, wherein the compound is represented by a structural formula VI.

or a pharmaceutically acceptable salt, solvate or hydrate thereof.

- 8. (Canceled)
- 9. (Canceled)

 (Currently amended). The compound of Claim 2, wherein the compound is represented by structural formula VIII,

$$\begin{array}{c|c}
 & R^{1} \\
\hline
 & R^{1} \\
\hline
 & A_{1} \\
\hline
 & COOR^{6}
\end{array}$$

VIII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein: A_1 and A_2 are respectively:

O and O.

CH2 and O,

CH2 and S.

O and S or

S and O:

m is: 1 or 2:

R1 is: C1-C2 alkyl; and

R3 is: hydrogen, halo or C1-C6 alkvl;

R6 and R9 are each independently; hydrogen or C1-C6 alkyl;

T is: a <u>bond or bond, -O-, -C(O) , S(O) S(O)2-, C(=CH2) , -C(=NOH) or -CH(OH)-;</u> and ring b is optionally substituted with one or more groups independently selected from:

hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, S(O)₂R⁹, C₁-C₆ alkyl, C₁-C₆ alkoxy and (CH₂)_nC₃-C₈ cycloalkyl.

 (Previously Presented). The compound of Claim 10, wherein the compound is represented by structural formula IX,

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

R1 is C1-C3 alkyl;

R3 is: hydrogen, halo or C1-C4 alkyl;

ring b is optionally substituted with one or more groups independently selected from the group consisting of: hydrogen, halo, haloalkyl, haloalkyloxy and C₁-C₆ alkyl.

12. (Previously Presented). The compound of Claim 11, wherein the compound is represented by structural formula X,

or a pharmaceutically acceptable salt, solvate or hydrate thereof.

 (Previously Presented). The compound of Claim 11, wherein the compound is represented by structural formula XI,

$$H_3C$$
 CH_3
 CH_3
 $COOH$

 (Withdrawn). The compound of Claim 2, wherein the compound is represented by structural formula XII,

$$\bigcap_{N = 1}^{R^1} \bigcap_{N = 1}^{R^3} \bigcap_{N = 1}^{R^3} \bigcap_{N = 1}^{COOR^6} \bigcap_{N = 1}^{R^4} \bigcap_{N = 1}^{R^6} \bigcap_{N = 1}$$

XII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein: A₁ and A₂ are respectively:

O and O,

CH2 and O.

CH2 and S.

O and S or

S and O:

S and O

m is: 1 or 2;

R1 is: C1-C3 alkyl; and

R3 is: hydrogen, halo or C1-C6 alkyl;

R4, R5, R6 and R9 are each independently: hydrogen or C1-C6 alkyl;

rings k and l are each optionally substituted with one or more groups independently selected from:

hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, S(O)₂R⁹, C₁-C₆ alkyl, C₁-C₆ alkoxy and (CH₂)_nC₃-C₈ cycloalkyl.

15. (Canceled)

 (Previously Presented). The compound of Claim 2, wherein the compound is represented by structural formula XIII,

$$Z \xrightarrow[O]{(CH_2)_m} A_2 \xrightarrow[R^4]{(R^3)_r} A_1 \xrightarrow[R^4]{COOR^6}$$

XIII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein m is 1, 2, 3, or 4.

17. (Canceled).

18. (Withdrawn). The compound of Claim 16, wherein the compound is represented by structural formula XV,

XV

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

T is: a bond, O or C(O);

R2 is: methyl, ethyl or cyclopropyl;

R3 is: methyl or ethyl; and

rings b and c are each optionally substituted with one or more substitutent independently selected from the group consisting of: hydrogen, Cl, Br, CF₃, OCF₃, N(CH₃)₂, S(O)₂CH₃, methyl, ethyl, isopropyl, methoxy and cyclopropyl.

 (Currently Amended). The compound of Claim 2, wherein the compound is represented by structural formula XVI,

XVI

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein Y is a branched alkyl-or C_k - C_k -eyeloalkyl.

20. (Canceled).

21. (Withdrawn). The compound of Claim 19, wherein the compound structural formula XVIII,

$$\begin{array}{c|c}
R^3 \\
\hline
COOH \\
R^{9a} R^{9b}
\end{array}$$

XVIII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

T is: a bond, O or C(O);

R3 is: methyl or ethyl;

 R^{9a} and R^{9b} are each independently hydrogen, methyl or ethyl, wherein at least one of R^{9a} and R^{9b} being methyl or ethyl;

- rings b and c are each optionally substituted with one or more substituent independently selected from the group consisting of: hydrogen, Cl, Br, CF₃, OCF₃, S(O)₂CH₃, N(CH₃)₂, methyl, ethyl, isopropyl, methoxy and cyclopropyl.
 - 22. (Canceled).
- (Withdrawn). The compound of Claim I, wherein the compound is a compound of formula XX,

$$Z = \bigcup_{Q = 1}^{R^1} \bigvee_{Q = 1}^{R^2} \bigvee_{Q = 1}^{(R^2)_r} \bigvee_{Q = 1}^{A_1} \bigvee_{Q = 1}^{Q} \bigvee_{Q = 1}^{Q} \bigvee_{Q = 1}^{A_1} \bigvee_{Q = 1}^{Q} \bigvee_{Q = 1}^$$

XX

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A₁ is: a bond, CH₂, O or S, and wherein A₁ and R⁴ or A₁ and R⁵ together being a 3- to 6membered carbocyclyl when A₁ is a carbon;

A2 is: O or S or CH2;

Q is: -C(O)OR⁶, or R^{6A};

Y is: a bond, C₁-C₆ alkyl or C₃-C₆ cycloalkyl;

Z is: a) arvl:

- a 5- to 10-membered heteroaryl wherein the heteroaryl containing at least one heteroatom selected from N, O or S.
- bi-aryl, wherein biaryl being defined as aryl substituted with another aryl or aryl substituted with heteroaryl, or
- d) bi-heteroaryl, wherein bi-heteroaryl being defined as heteroaryl substituted with another heteroaryl, or heteroaryl substituted with aryl, and wherein aryl, heteroaryl, bi-aryl and bi-heteroaryl being optionally substituted with one or more groups independently selected from R⁷;

n is: 1, 2, 3, 4, 5 or 6

p is: 1 or 2;

r is: 1, 2, 3, or 4;

R1 and R2 are each independently:

hydrogen,

haloalkyl,

C1-C6 alkyl,

(CH₂)_nC₃-C₈ cycloalkyl, or

 R^1 and R^2 form a 4- to 8-membered nonaromatic carbocyclic ring; and wherein at least one of R^1 and R^2 is alkyl or cycloalkyl, and;

R3 is: hydrogen,

nitro,

cyano,

hydroxyl,

halo.

haloalkyl,

haloalkyloxy,

aryloxy,

C₁-C₆ alkyl,

C1-C6 alkoxy or

C₃-C₈ cycloalkyl;

R4 and R5 are each independently; hydrogen or C1-C6 alkyl;

R6 is: hydrogen, C1-C6 alkyl or aminoalkyl;

R^{6A} is: carboxamide, sulfonamide, acylsulfonamide, tetrazole,

R7 is: hydrogen,

oxo,

nitro,

cyano,

hydroxyl,

halo,

haloalkyl,

haloalkyloxy,

aryloxy,

arylalkyl,

aminoalkyl,

C₁-C₆ alkyl,

C₁-C₆ alkoxy,

(CH₂)_nC₃-C₈ cycloalkyl,

 $C(O)R^9$,

C(O)OR⁹, C(=NOR⁸)R⁹, CR⁸(OH)R⁹, C[=C(R⁸)₂]R⁹, OR⁹, SR⁹ or S(O)₀R⁹;

R8 is: hydrogen or C1-C6 alkyl; and

 R^9 is: hydrogen, C_1 - C_6 alkyl, C_3 - C_8 cycloalkyl, aryl, heteroaryl or

heterocyclyl.

wherein alkyl, cycloalkyl, aryl, heteroaryl or heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of:

hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1 - C_6 alkyl, C_1 - C_6 alkoxy and C_3 - C_8 cycloalkyl.

- 24. (Canceled).
- 25. (Canceled).
- (Withdrawn). The compound of Claim 23, wherein the compound is a compound of structural formula XXII,

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

T is: a bond, -O- or -C(O)-;

R1 is: methyl, ethyl or cyclopropyl;

R3 is: methyl or ethyl; and

rings b and c are each optionally substituted with one or more substituent independently selected from the group consisting of: hydrogen, Cl, Br, CF₃, OCF₃, S(O)₂CH₃, N(CH₃)₂, methyl, ethyl, isopropyl, methoxy and cyclopropyl.

 (Withdrawn). The compound of Claim 1, wherein the compound is a compound of structural formula XXIII,

$$\begin{array}{c}
R^1 \\
b \\
COOR^4
\end{array}$$

$$\begin{array}{c}
R^3 \\
A_1 \\
COOR^4
\end{array}$$

$$\begin{array}{c}
COOR^4 \\
XXIII
\end{array}$$

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A₁ and A₂ are respectively:

O and O,

CH2 and O,

CH2 and S.

O and S or

S and O:

m is: 1, 2, 3 or 4;

R1 is: C1-C3 alkyl; and

R3 is: hydrogen, halo or C1-C6 alkyl;

R6 and R9 are each independently: hydrogen or C1-C6 alkyl;

 $T\ is:\ a\ bond, -O-, -C(O)-, -S(O)-S(O)_2-, -C(=CH_2)-, -C(=NOH)-\ o\ -CH(OH)-;\ and\ rings\ b\ and\ c\ are\ each\ optionally\ substituted\ with\ one\ or\ more\ groups\ independently\ selected\ from:$

hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, $S(O)_2R^9$, $C_{1^*}C_6$ alkyl, $C_{1^*}C_6$ alkoxy and $(CH_2)_nC_3-C_8$ cycloalkyl.

- 28. (Canceled).
- (Currently Amended). A compound of Claim 1 selected from the group consisting of:

	Structure	Name
1	H ₂ C CH ₃ OH	3 (4 [3 (2 Benzoyl 4 ethyl- phenoxy) butoxy] 2- methyl phenyl) propionie acid
2	H ₃ C OH ₃	[4 [3 (2-Benzoyl 4 ethyl- phenoxy) butoxy] 2- methyl-phenoxy] acetic acid
3	H ₃ C O CH ₃ O O O O O O O O O O O O O O O O O O O	14 [3 (2 Benzoyl 4 ethyl- phenoxy)-butoxy]-2- methyl-phenylsulfunyl)- acetic acid

	Structure	Name
4	H ₃ C CH ₃ OH	[44[3-(2-Benzoyl-4-ethyl- phenoxy)-butoxy]-2- methyl-phenylsulfanyl}- acetic-acid
5	H ₃ C S CH ₃ O O O O O O O O O O O O O O O O O O O	[4-[3-(2-Benzoyl-4-ethyl- phenoxy)-butylsulfanyl]-2- methyl-phenoxy]-acetic acid
6	H ₃ C S OH	3 (1 [3 (2-Benzoyl 4 ethyl- phenoxy)-butylsulfanyl] 2- methyl-phenyl]-propionie acid
7	H ₃ C CH ₃ O CH ₃ O H ₄ O CH ₃ O C	2-[4-[3-(2-Benzoyl-4-ethyl- phenoxy) butoxy] 2- methyl-phenoxy] 2-methyl- propionic-acid
8	H,C O O O O O O O O O O O O O O O O O O O	14 [3 (2 Benzoyl 4 ethyl- phenoxy) butoxy] phenoxy) acetic acid

	Structure	Name
10	H ₃ C CH ₃ OH Chiral	3-[4-[3-(2-Benzoyl-4-isopropyl-phenoxy)-butoxy]-2-methyl-phenyll-propionic-acid 3-[4-[3-(2-Benzoyl-4-eyelopropyl-phenoxy)-
	CH ₃ OH	butoxy] 2 methyl phenyl}- propionie acid
11	F CH ₃ OH	3 - [4 [3 (2 Benzoyl 4 trifluoromethyl-phenoxy)- butoxy] 2 methyl-phenyl}- propionic acid
12	CI—CH ₃	3- [4-[3-(2-Benzoyl-4- ehloro-phenoxy)-butoxy]-2- methyl-phenyl)-propionie acid
13	CI—CH ₃	3-(4-[3-(2-Benzoyl-4- ehloro-phenoxy)-butoxy]-2- methyl-phenyl]-propionie aeid

	Structure	Name
14	Chiral Chiral CH ₃ OH	3-[4-[3-(2-Benzoyl-4- methoxy-phenoxy)- butoxy] 2-methyl-phenyl}- propionic-acid
15	Chiral Chiral	3 [4 [3 (2-Benzoyl 4- fluoro-phenoxy)-butoxy] 2- methyl-phenyl}-propionie acid
16	H ₃ C CH ₃ OH	3-[4-[3-(2-Benzoyl-4- isopropyl-phenoxy)- butoxy]-2-methyl-phenyl}- propionie-acid
17	CI-CH ₃ O-CH ₃ OH	[4 [3 (2-Benzoyl 4-chloro- phenoxy)-butoxy] 2- methyl-phenylsulfanyl)- acetic-acid
18	H ₃ C OH OH OH	3 (4 13 [4 Ethyl 2- (hydroxy-phenyl-methyl)- phenoxy]-butoxyl-2- methyl-phenyl)-propionic acid

	Structure	<u>Name</u>
19	H ₂ C OH OH	3 (4 [3 [4 Ethyl 2 (hydroxyimino phenyl methyl) phenoxy] butoxy] 2 methyl phenyl) propionic acid
20	H ₃ C CH ₃ OH	3 (4-[3-[4-Ethyl-2- (methoxyimino phenyl- methyl) phenoxy] butoxy]- 2-methyl-phenyl) propionie acid
21	H ₂ C CH ₃ Christ	3-[4-[3-(4-Isopropyl-2- phonoxy phonoxy) butoxy] 2-methyl phonyl} propionic acid
22	H ₂ C Chral	[4 [3 (4 Isopropyl 2 phenoxy phenoxy) butoxy] 2 methyl- phenykalfanyl] acetic acid
23	H ₃ C CH ₃ O CH ₃ O O O O O O O O O O O O O O O O O O O	3 14 [3 (4 Ethyl 2- isobutyryl phenoxy)- butoxy] 2 methyl phenyl}- propionic acid
24	H ₃ C CH ₃ OH	3 (4 [3 (2 Cyclopropanecarbonyl 4- ethyl phenoxy) butoxy] 2- methyl phenyl propionic acid

	Structure	<u>Name</u>
25	H ₅ C CH ₃ O CH ₃	3-[4-[3-(2- Cyclopropunecarbonyl 4- ethyl-phenoxy) butoxy] 2- methyl-phenyl -propionic acid
26	H ₃ C OH ₃	3 (4 [3 (2- Cyclopentanecarbonyl 4- ethyl-phenoxy)-butoxy]-2- methyl-phenyl -propionie acid
27	H ₃ C CH ₃	2-{4-{3-(4-Ethyl-2- isobutyryl-phenoxy)- butoxy]-phenoxy}-2- methyl-propionic acid
28	H ₃ C CH ₃ OH	2-[4-[3-(2- Cyclopropaneearbonyl-4- ethyl-phenoxy) butoxy] phenoxy] 2-methyl- propionic acid
29	H ₃ C CH ₃ OCH ₃ OCH ₃	3-{4-[3-(3-Benzoyl-5-ethyl- pyridin-2-yloxy)-butoxy]-2- methyl-phenyl}-propionic acid
30	H ₃ C CH ₃ OCH ₃ OCH	{4-{3-(3-Benzoyl-5-ethyl- pyridin-2-yloxy)-butoxy]-2- methyl-phenylsulfanyl}- acetic acid

	Structure	<u>Name</u>
31	Chiral	3-{4-[3-(3-Benzoyl-5-
		chloro-pyridin-2-yloxy)-
	CH³	butoxy]-2-methyl-phenyl}-
	CI-CI-CI-CI-CI-CI-CI-CI-CI-CI-CI-CI-CI-C	propionic acid
	Ċн _з	
22	OH Chiral	(452 (2 P) 15 11
32	Cimai	{4-[3-(3-Benzoyl-5-chloro-
		pyridin-2-yloxy)-butoxy]-2-
	CH ₃	methyl-phenylsulfanyl}-
	S. CH ₃	acetic acid
	ОН	
33	Chiral	3-{4-[3-(3-Benzoyl-5-
		trifluoromethyl-pyridin-2-
		yloxy)-butoxy]-2-methyl-
		phenyl}-propionic acid
	F CH ₃	
34	Chiral	{4-[3-(3-Benzoyl-5-
34		trifluoromethyl-pyridin-2-
	Ĭ.	yloxy)-butoxy]-2-methyl-
	FCH ₃	phenylsulfanyl}-acetic acid
		phonyisunanyiy-acciic acid
	H _s C	
35	Chiral	3-{4-[3-(5-Chloro-3-
	<u> </u>	phenoxy-pyridin-2-yloxy)-
	/ _ Сн₃	butoxy]-2-methyl-phenyl}-
	ci—()—o	propionic acid
	CH ₃	
	ОН	

	Structure	Name
36	Chiral	3-{4-[3-(5-Chloro-3-
	<u> </u>	phenoxy-pyridin-2-yloxy)-
	,CH ₃	butoxy]-2-ethyl-phenyl}-
		propionic acid
	CH,	
37	Chiral	{4-[3-(5-Chloro-3-
"	_>	phenoxy-pyridin-2-yloxy)-
	, CH ₃	butoxy]-2-methyl-
		phenylsulfanyl}-acetic acid
	" H ₃ C — S— S	priesty to arranger, we come work
	о́н	
38	F CH ₃ Chiral	3-{2-Methyl-4-[3-(3-
		phenoxy-5-trifluoromethyl-
	O CH ₃ OH	pyridin-2-yloxy)-butoxy]-
		phenyl}-propionic acid
39	Chiral	3-{2-Ethyl-4-[3-(3-
	_ \'	phenoxy-5-trifluoromethyl-
	Fócн₃	pyridin-2-yloxy)-butoxy]-
		phenyl}-propionic acid
	F N H ₃ C OH	
40	Chiral	3-{2-Ethyl-4-[3-(3-
		phenoxy-5-trifluoromethyl-
	,cH ₃	pyridin-2-yloxy)-butoxy]-
		phenyl}-propionic acid
	F N H ₃ C	
	'он	
41	5	3-{2-Methyl-4-[3-(3-
	, Ł , OH	phenoxy-5-trifluoromethyl-
	F, CH3	pyridin-2-yloxy)-propoxy]-
		phenyl}-propionic acid
		(trifluoroacetic acid salt)
	, он	

	Structure	Name
42	F-FOH CI-CH ₅	3-{4-[3-(5-Chloro-3- phenoxy-pyridin-2-yloxy)- propoxy]-2-methyl- phenyl}-propionic acid
43	CI——N OH	3-{4-[2-(5-Chloro-3- phenoxy-pyridin-2- ylamino)-ethoxy]-2-methyl- phenyl}-propionic acid
44	H ₃ C CH ₃ OH	3-{4-[3-(3-Benzoyl-5-ethyl- pyridin-2-yloxy)-propoxy]- 2-methyl-phenyl}-propionic acid
45	Chiral Chiral Chiral	3-{2-Methyl-4-[3-(6- methyl-2-phenoxy-pyridin- 3-yloxy)-butoxy]-phenyl}- propionic acid
46	H ₅ C CH ₅ OCH ₅	3 (4 [3 (5 Ethyl biphenyl 2 yloxy) butoxy] 2 methyl- phenyl} propionic acid

	Structure	<u>Name</u>
47	Chiral Chiral OH	3 [4-[3 (4-Ethyl-2-oxazol- 2 yl-phenoxy) butoxy] 2- methyl-phenyl}-propionie acid
48	H ₂ C CH ₃ Christ	3 [4 [3 (4 Ethyl 2 thiazol 4 yl phenoxy) butoxy] 2-methyl phenyl) propionie aeid
49	Chiral H ₂ C CH ₃ OH	3-{4-[3-(4-Ethyl-2-pyridin- 2-yl-phenoxy)-butoxy]-2- methyl-phenyl}-propionic acid
50	H ₃ C Chiral	{4-[3-(4-Ethyl-2-pyridin-2- yl-phenoxy)-butoxy]-2- methyl-phenylsulfanyl}- acetic acid
51	CH ₃ Chiral OH	3-{2-Ethyl-4-[3-(4-ethyl-2- pyridin-2-yl-phenoxy)- butoxy]-phenyl}-propionic acid
52	Chiral Chiral Chiral	3-{4-[3-(4-Chloro-2- pyridin-2-yl-phenoxy)- butoxy]-2-methyl-phenyl}- propionic acid

	Structure	Name
53	F CH ₃ Chiral	3-{2-Methyl-4-[3-(2- pyridin-2-yl-4- trifluoromethyl-phenoxy)- butoxy]-phenyl}-propionic acid
54	F F O O O O O O O O O O O O O O O O O O	actu 3-{2-Ethyl-4-[3-(2-pyridin- 2-yl-4-trifluoromethyl- phenoxy)-butoxy]-phenyl}- propionic acid
55	H ₂ C CH ₃ OH	3-{4-{3-(4-Ethyl-2-pyridin- 3-yl-phenoxy)-butoxy]-2- methyl-phenyl}-propionic acid
56	CI—CH ₃ OH	3-{4-[3-(4-Chloro-2- pyridin-3-yl-phenoxy)- butoxy]-2-methyl-phenyl}- propionic acid
57	H ₂ C CH ₂ OH	3-{4-{3-(4-Ethyl-2-pyridin- 4-yl-phenoxy)-butoxy]-2- methyl-phenyl}-propionic acid
58	F CH ₃ Chiral Chiral OH	3-(2-Methyl-4-(3-(2- pyridin-4-yl-4- trifluoromethyl-phenoxy)- butoxy]-phenyl}-propionic acid

	Structure	<u>Name</u>
59	F H ₃ C Chiral	3-{2-Ethyl-4-[3-(2-pyridin- 4-yl-4-trifluoromethyl- phenoxy)-butoxy]-phenyl}- propionic acid
60	CH ₃ Chiral	3 [4 [3 (2- Benzo[d]isoxazol 3 yl 4- chloro phenoxy) butoxy] 2- methyl phenyl) propionie aeid
61	H ₂ C OH OH	3-(4-[3-(2-Benzoyl-4-ethyl- phenoxy)-butoxy]-2- methyl-phenyl -propionic acid
62	H ₃ C CH ₃ O-CH ₃ OH	14 [3 (2 Benzoyl 4 ethyl- phenoxy) butoxy] 2- methyl-phenoxy] acetic acid
63	H ₂ C CH ₃ OCH ₃	[4 [3 (2 Benzoyl 4 ethyl- phenoxy) butoxy] 2- methyl-phenylsulfanyl)- acetic acid
64	H ₃ C CH ₃ O CH ₃	[4 [3 (2-Benzoyl 4 ethyl- phenoxy) butoxy] 2- methyl-phenylsulfanyl]- acetic acid

	Structure	Name
65	H ₃ C S CH ₃ O O O O O O O O O O O O O O O O O O O	[4-[3-(2-Benzoyl-4-ethyl- phenoxy)-butylsulfanyl] 2- methyl-phenoxy]-acetic acid
66	H ₃ C S OH ₃	3 [4 [3 (2 Benzoyl 4 ethyl- phenoxy)-butylsulfanyl] 2- methyl-phenyl]-propionie acid
67	H ₃ C CH ₃ O CH ₃ O O O O O O O O O O O O O O O O O O O	2 (4 [3 (2 Benzoyl 4 ethyl- phenoxy) butoxy] 2- methyl-phenoxy] 2 methyl- propionic acid
68	H ₃ C O O O O O O O O O O O O O O O O O O O	14 [3 (2 Benzoyl 4 ethyl- phenoxy) butoxy]- phenoxy] acetic acid
69	H ₃ C CH ₃ OCH ₃	3-{4-[3-(2-Benzeyl-4- isopropyl-phenoxy)- butoxy]-2-methyl-phenyll- propionic-acid

	Structure	<u>Name</u>
70	Chiral	3-[4-[3-(2-Benzoyl-4-
		eyclopropyl-phenoxy)
	CH ₃	butoxy]-2-methyl-phenyl}-
		propionic acid
	CH ₃	
71		3-14-[3-(2-Benzoyl-4-
'		trifluoromethyl-phenoxy)-
		butoxy] 2 methyl phenyl)
	CH ₃	propionie acid
	F CH ₃	
	OH OH	
72	/=\	3 {4 [3 (2 Benzoyl 4
		ehloro phenoxy) butoxy] 2
	O	methyl-phenyl}-propionie
	c c	acid
	CH, OH	
73		3-{4-[3-(2-Benzoyl-4-
"		chloro-phenoxy) butoxy] 2-
)=o	methyl-phenyl}-propionic
	ci—Co	aeid
	CH, CO	
	он	
74	Chiral	3 [4 [3 (2 Benzoyl 4
		methoxy phenoxy)
	CH ₃	butoxy]-2-methyl-phenyl}-
	H ₃ C-O	propionic acid
	ČH ₃ OH	
	L	

	Structure	Name
	Chiral	
75	Chiral	3-[4-[3-(2-Benzoyl-4-
		fluoro-phenoxy) butoxy] 2-
	>	methyl-phenyl}-propionic
	CH ₃	acid
	CH ₃	
76	Chiral	3-{4-{3-(2-Benzoyl-4-
70		
		isopropyl-phenoxy)-
	H ₃ C CH ₃	butoxy] 2 methyl phenyl}-
	H,C O	propionie-acid
	CH ₃	
77	Chiral	(4.52.(2.D. 1.4
++	Cilia	(4 [3 (2 Benzoyl 4
		isopropyl-phenoxy)-
	CH ₃	butoxy]-2-methyl-
	H ₃ C S S	phenylsulfanyl)-acetic-acid
	H ₃ Ć CH ₃ OH	
78		(4-[3-(2-Benzoyl-4-chloro-
		phenoxy)-butoxy]-2-
	<u> </u>	methyl-phenylsulfunyl}-
	CI CH ₃ Q	acetic acid
	CH ₂ O-S OH	
	G13	A (1 (2 51 m) 1 A
79		3 (4 {3 [4 Ethyl 2
		(hydroxy-phenyl-methyl)-
	— ОН СН3	phenoxy] butoxy) 2
	H ₃ C CH ₃	methyl-phenyl) propionie
	CH0-	acid
	`он	

	Structure	<u>Name</u>
80	N-OH CH ₃	3-(4-[3-[4-Ethyl-2- (hydroxyimino-phenyl- methyl)-phenoxy]-butoxy}- 2-methyl-phenyl)-propionie aeid
81	H ₃ C CH ₃ OCH ₃ OH	3 (4 [3 [4 Ethyl 2- (methoxyimino phenyl- methyl) phenoxy] butoxy}- 2 methyl phenyl) propionie acid
82	H ₃ C CH ₃ Chiral	3-[4-[3-(4-Isopropyl-2- phenoxy-phenoxy-) butoxyl-2-methyl-phenyll- propionic-acid
83	H ₂ C OH OH	14 [3 (4 Isopropyl 2- phenoxy phenoxy) butoxy] 2 methyl- phenykulfanyl] acetic acid
84	H ₃ C, CH ₃	3-{4-[3-(4-Ethyl-2- isobutyryl-phenoxy)- butoxy]-2-methyl-phenyl}- propionic-acid
85	H ₂ C OH ₃	3-[4-[3-(2- Cyclopropanecarbonyl-4- ethyl-phenoxy)-butoxy]-2- methyl-phenoyl]-propionic acid

	Structure	Name
86	\triangleleft	3-[4-[3-(2-
)=o	Cyclopropanecarbonyl-4-
	H ₃ C CH ₃	ethyl-phenoxy) butoxy]-2-
	CH ₃	methyl-phenyl -propionic
	ОН	acid
87	\bigcap	3-{4-{3-(2-
	\ \ \	Cyclopentanecarbonyl-4
	н,с,	ethyl-phenoxy) butoxy]-2-
		methyl-phenyl}-propionie
	H,c _0	acid
	ОН	
88	H₃C ✓ CH₃	2-{4-[3-(4-Ethyl-2-
	H.G - 9	isobutyryl-phenoxy)
	~	butoxy] phenoxy) 2
	CH ₃ H ₃ C CH ₃	methyl-propionic acid
89	◁	2 (4 [3 (2
) _0	Cyclopropanocarbonyl 4
	# ₅ c	ethyl-phenoxy)-butoxy]-
		phenoxy}-2-methyl-
	CH ₃ H ₃ C CH ₃ OH	propionie acid
90	H₃C /=N ,CH₃	3-{4-[3-(3-Benzoyl-5-ethyl-
		pyridin-2-yloxy)-butoxy]-2-
	OH OH	methyl-phenyl}-propionic
		acid
91	/=\	{4-[3-(3-Benzoyl-5-ethyl-
		pyridin-2-yloxy)-butoxy]-2-
	H.C. CH ₃	methyl-phenylsulfanyl}-
	H ₃ C	acetic acid
	CH ₃	
	ОН	

	Character and Ch	N
	Structure	Name
92	Chiral	3-{4-[3-(3-Benzoyl-5-
		chloro-pyridin-2-yloxy)-
	Сн,	butoxy]-2-methyl-phenyl}-
		propionic acid
	CH ₃	
	он	
93	Chiral	{4-[3-(3-Benzoyl-5-chloro-
		pyridin-2-yloxy)-butoxy]-2-
	,cH ₁	methyl-phenylsulfanyl}-
	cı—()	acetic acid
	CH ₃	
	он	
94	Chiral	3-{4-[3-(3-Benzoyl-5-
		trifluoromethyl-pyridin-2-
	, сн,	yloxy)-butoxy]-2-methyl-
		phenyl}-propionic acid
	F _N CH3 OH	
95	Chiral	{4-[3-(3-Benzoyl-5-
		trifluoromethyl-pyridin-2-
		yloxy)-butoxy]-2-methyl-
	CH ₃	phenylsulfanyl}-acetic acid
	F H ₁ C OH	, , , ,
06	Chiral	2 (4 52 (5 Chloro 2
96		3-{4-[3-(5-Chloro-3-
		phenoxy-pyridin-2-yloxy)-
	Сн³	butoxy]-2-methyl-phenyl}-
		propionic acid
	t _H 3 OH	
		l .

	Structure	Name
97	Chiral	3-{4-[3-(5-Chloro-3-
	>= /	phenoxy-pyridin-2-yloxy)-
	CH³	butoxy]-2-ethyl-phenyl}-
		propionic acid
	N CH3 OH	
98	Chiral	{4-[3-(5-Chloro-3-
	>= /	phenoxy-pyridin-2-yloxy)-
	, CH₃	butoxy]-2-methyl-
	CI-V _N O _{H₃C} OO _O OO	phenylsulfanyl}-acetic acid
99	F /=N CH ₃ Chiral	3-{2-Methyl-4-[3-(3-
"		phenoxy-5-trifluoromethyl-
	CH ₃	pyridin-2-yloxy)-butoxy]-
		phenyl}-propionic acid
100	Chiral	
100	()	3-{2-Ethyl-4-[3-(3-
		phenoxy-5-trifluoromethyl-
	сн,	pyridin-2-yloxy)-butoxy]- phenyl}-propionic acid
	F N HLC	phenyry-proprome acid
	, он	
101	Chiral	3-{2-Ethyl-4-[3-(3-
		phenoxy-5-trifluoromethyl-
	F_CH ₃	pyridin-2-yloxy)-butoxy]-
	F N H,C	phenyl}-propionic acid
	ОН	
102		3-{2-Methyl-4-[3-(3-
	FF OH	phenoxy-5-trifluoromethyl-
	F, CH ₃	pyridin-2-yloxy)-propoxy]-
	F N O O O	phenyl}-propionic acid
	L C	(trifluoroacetic acid salt)
		l .

	Structure	Name
103	FJ 0	3-{4-[3-(5-Chloro-3-
	FOH	phenoxy-pyridin-2-yloxy)-
		propoxy]-2-methyl-
	,CH,	phenyl}-propionic acid
	CI-(N)	
	ОН	
104		3-{4-[2-(5-Chloro-3-
		phenoxy-pyridin-2-
	,cн _а	ylamino)-ethoxy]-2-methyl-
	CI—N	phenyl}-propionic acid
105	H ₂ C, /=N, ,CH ₃	3-{4-[3-(3-Benzoyl-5-ethyl-
103		pyridin-2-yloxy)-propoxy]-
		2-methyl-phenyl}-propionic
	ОН	acid
	y	
106	Chiral	3-{2-Methyl-4-[3-(6-
	,	methyl-2-phenoxy-pyridin-
	H C CH ₃	3-yloxy)-butoxy]-phenyl}-
	H.C.	propionic acid
	он —	
107		3 {4 [3 (5 Ethyl biphenyl
	<u>"</u>	2 yloxy) butoxy] 2 methyl-
	H ₃ C CH ₃	phenyl)-propionie acid
	CH ₃	
108	OH Chiral	3 4 3 (4 Ethyl 2 oxazol-
108		
	H,C,	2-yl-phenoxy) butoxy] 2-
	OH OH	methyl-phenyl) propionie
	ČH ₃	acita

	Structure	Name
109	H ₃ C Chiral	3-[4-[3-(4-Ethyl-2-thiazol-
		4-yl-phenoxy) butoxy] 2-
	CH ₃	methyl-phenyl}-propionic
	-5	aeid
110	Chiral	3-{4-[3-(4-Ethyl-2-pyridin-
) N	2-yl-phenoxy)-butoxy]-2-
	H ₃ C CH ₃	methyl-phenyl}-propionic
	H ₂ C OH	acid
111	Chiral	{4-[3-(4-Ethyl-2-pyridin-2-
	CH3 8	yl-phenoxy)-butoxy]-2-
	H ₃ C S OH	methyl-phenylsulfanyl}-
	H ₀ C	acetic acid
112	CH ₂ Chiral	3-{2-Ethyl-4-[3-(4-ethyl-2-
112	CH, Sima	pyridin-2-yl-phenoxy)-
	ОН	butoxy]-phenyl}-propionic
		acid
	Ĉн₃	acid
113	Chiral	3-{4-[3-(4-Chloro-2-
	>=n'	pyridin-2-yl-phenoxy)-
	CI—CI—O	butoxy]-2-methyl-phenyl}-
	H _s c O	propionic acid
	он	
114	F, CH ₃ Chiral	3-{2-Methyl-4-[3-(2-
	F OH	pyridin-2-yl-4-
		trifluoromethyl-phenoxy)-
	CH ₃	butoxy]-phenyl}-propionic
	110	acid
115	F H ₃ C Chiral	3-{2-Ethyl-4-[3-(2-pyridin-
	F	2-yl-4-trifluoromethyl-
	~ ° ~ °	phenoxy)-butoxy]-phenyl}-
	Сн _з	propionic acid

	Structure	Name
116	H ₃ C Chral	3-{4-[3-(4-Ethyl-2-pyridin- 3-yl-phenoxy)-butoxy]-2- methyl-phenyl}-propionic acid
117	CI—CH ₅ CH ₅ OH	3-{4-[3-(4-Chloro-2- pyridin-3-yl-phenoxy)- butoxy]-2-methyl-phenyl}- propionic acid
118	H ₂ C CH ₂ OH	3-{4-[3-(4-Ethyl-2-pyridin- 4-yl-phenoxy)-butoxy]-2- methyl-phenyl}-propionic acid
119	F Christ	3-{2-Methyl-4-[3-(2- pyridin-4-yl-4- trifluoromethyl-phenoxy)- butoxy]-phenyl}-propionic acid
120	F CH ₃	3-{2-Ethyl-4-[3-(2-pyridin- 4-yl-4-trifluoromethyl- phenoxy)-butoxy]-phenyl}- propionic acid
121	Cl Chiral OH	3-[4-[3-(2- Benzo[d]isoxazol 3 yl 4- chloro phenoxy) butoxy] 2- methyl phenyl} propionie aeid

122	
phenoxy phenoxy) butoxy] 2-methyl phenylsulfanyl) acetic- Cha Chrol (R) 14 [3 (2-benzoyl 4-thyl phenylsulfany) acetic-acid (R) [4 [3 (2-benzoyl 4-thyl phenylsulfany) acetic-acid (R) [4 [3 (2-benzoyl 4-thyl phenylsulfany)] acetic-acid (R) [4 [3 (2-benzoyl 4-thyl phenylsulfany]] acetic-acid (R) [4 [3 (2-benzoyl 4-thyl phenylsulfany]] acetic-acid (R) [4 [3 (2-benzoyl 4-thyl phenylsulfany]] acetic-acid	
H ₃ C OH CH ₃ Duttoxy] 2-methyl-phenylsulfanyl}-acetic- CH ₃ Chiral (R) 14 [3 (2-benzoyl-4-methyl-phenylsulfany)] acetic-acid Ch ₃ Christ (R) [4 [3 (2-benzoyl-4-trifluoromethoxy-phene-butoxy] 2-methyl-phenylsulfany) acetic-acid (R) [4 [3 (2-benzoyl-4-trifluoromethoxy-phene-butoxy] 2-methyl-phenylsulfanyl] acetic-acid (A) [4 [3 (2-benzoyl-4-trifluoromethoxy-phene-butoxy] 2-methyl-phenylsulfanyl] acetic-acid	
H ₃ C OH CH ₃ OH Chiral (R) 4 [3 (2 benzoyl 4 urifluoromethoxy phenolsulfanyl) acetic-acid CH ₃ Chiral (R) 4 [3 (2 benzoyl 4 urifluoromethoxy phenolsulfany acetic-acid (R) 4 [3 (2 benzoyl 4 urifluoromethoxy phenolsulfany) phenolsulfanyl) acetic-acid (R) 4 [3 (2 benzoyl 4 urifluoromethoxy phenolsulfanyl) acetic-acid (R) 4 [3 (2 benzoyl 4 urifluoromethoxy phenolsulfanyl) acetic-acid (R) 4 [3 (2 benzoyl 4 urifluoromethoxy phenolsulfanyl) acetic-acid	
123 Chiral Chiral Chiral (R) 14 [3 (2-benzoyl 4-methyl-phenoxy) buttox 2-methyl-phenoxy) buttox 2-methyl-phenoxy) buttox 2-methyl-phenoxy) buttox 4-methyl-phenoxy) buttox 2-methyl-phenoxy) buttox 2-methyl-phenoxy phenobuttoxy-phenobuttoxy-phenobuttoxy-phenobuttoxy-phenobuttoxy-phenobuttoxy-phenoyl-acetic-acid (R) [4 [3 (2-benzoyl 4-triphenylsulfanyl) acetic-acid (R) [4 [3 (2-benzoyl 4-triphenylsulfanyl) acetic-acid	
123 Christ (R) 14 [3 (2-benzoyl-4 methyl-phenoxy) buttox 2-methyl-phenoxy) buttox 2-methyl-phenoxy) buttox 2-methyl-phenylsulfany acetic-acid (R) [4 [3 (2-benzoyl-4 trifluoromethoxy-phenobuttoxy] 2-methyl-phenylsulfanyl) acetic-acid (H) [4 [3 (2-benzoyl-4 trifluoromethoxy-phenobuttoxy] 2-methyl-phenylsulfanyl) acetic-acid (H) [4 [3 (2-benzoyl-4 ethyphenoxy) bexyloxy] 2-methyl-phenylsulfanyl) acetic-acid	eid
H ₃ C CH ₃ OH CH ₃ CH ₃ CH ₄ Chenzoyl 4 trifluoromethoxy-pheno butoxy] 2-methyl-phenylsulfany acetic-acid (R) [4 [3 (2-benzoyl 4 trifluoromethoxy-pheno butoxy] 2-methyl- phenylsulfanyl] acetic-acid (H) [4 [3 (2-benzoyl 4-ethyphenoxy) hexyloxy] 2-methyl-phenylsulfanyl] acetic-acid	
methyl phenoxy) butex 2-methyl phenylsulfany aeetie-aeid Chral (R) [4 [3 (2-benzoyl 4 trifluoromethoxy-pheno butoxy] 2-methyl- phenylsulfanyl) aeetie- (H) [4 [3 (2-benzoyl 4 trifluoromethoxy-pheno butoxy] 2-methyl- phenylsulfanyl) aeetie- aeetie-aeid	
2-methyl-phenylsulfany acetic-acid 124 Chiral (R) [4 [3 (2-benzoyl-4 trifluoromethoxy-phenobutoxy]-2-methyl-phenylsulfanyl) acetic-acid (R) [4 [3 (2-benzoyl-4 trifluoromethoxy-phenobutoxy]-2-methyl-phenylsulfanyl) acetic-acid	
H ₂ C O CH ₃ O CH ₃ Proposition of CH ₃ H ₂ C O CH ₃ Proposition of CH ₃ H ₃ C O CH ₃ Chiral (R) [4 [3 (2-benzoyl-4 trifluoromethoxy-phenobutoxy] 2-methyl-phenylsulfanyl] acetical decical	-
CH ₃ OH Chral (R) [4-[3-(2-benzoyl-4 triffluoromethoxy-phence butoxy]-2-methyl-phenylsulfunyl] acetic-acid (R) [4-[3-(2-benzoyl-4-ethyphenylsulfunyl] acetic-acid	H-
124 Chiral (R) [4 [3 (2-benzoyl-4 trifluoromethoxy-phenology-phe	
The content of the co	
FOCH3 butoxy] 2-methyl-phenylsulfunyl) acetic-to-to-to-to-to-to-to-to-to-to-to-to-to-	
H ₂ C OH ₃ OH Butoxy J 2 methyl phenylsulfunyl) acetic of the phenoxy hexyloxyl 2 methyl phenoxy) hexyloxyl 2 methyl phenoxy) hexyloxyl 2 methyl phenylsulfunyl) acetic acid	xy) -
Phenylsulfanyl) acetical phenylsulfanyl) acetical phenylsulfanyl) acetical [4 [3 (2 benzoyl 4 ethyphenowy) hexyloxy] 2 methyl phenylsulfanyl) aceticacid	
125 CH ₃ OH (4 [3 (2 benzoyl 4 ethy phenoxy) hexyloxy] 2-methyl phenylsulfanyll acetic acid	eid
125 (4 [3 (2 benzoyl 4 ethy phenoxy) hexyloxy] 2 methyl phenylsulfanyl) acetic-acid	
phenoxy) bexyloxy] 2- methyl phenylsulfanyl) acetic acid	
H ₃ C CH ₃ methyl-phenylsulfanyl} acetic-acid	<u> </u> -
H _s C O O O O O O O O O O O O O O O O O O O	
ОН	-
Ċн _з	
126 3-{4-{3 (2-benzoyl-4-et	n yl-
phenoxy) hexyloxy] 2	
CH ₃ methyl phenyl) propior	ie
H,C acid	
OH	
CH ₃	

	Structure	Name
127	Chiral	(R) 3-{4-[3-(4-ethyl-2-
		phenoxy-phenoxy)-
	H ₃ C, /=\ CH ₃	butoxy] 2-methyl-phenyl}-
		propionic acid
	CH ₃ OH	A.A.
128	Chiral	(R) 3 (4-13 [4 ethyl-2 (1-
		phenyl-vinyl) phenoxy]-
	H ₃ C CH ₂ CH ₃	butoxy} 2 methyl phenyl)
		propionic-acid
	ĊН ₃ ОН	
129	Chiral	(R) 3 (4 {3 [4 ethyl 2 (1
	CH ₃	methyl 1 phenyl ethyl)
	H ₃ C CH ₃ CH ₃	phenoxy] butoxy) 2
		methyl-phenyl) propionie
	CH ₃ OH	acid
130	Chiral	(R) 3 (4 [3 (2 benzoyl-4
		methyl-phenoxy)-butoxy]-
		2-methyl-phenyl)-propionie
	H ₃ C	acid
	CH ₃ OH	
140	Chiral	(R)-3-(4-[3-[4-ethyl-2-(1-
140		phenyl ethyl) phenoxy]
	CHCH_	butoxy\ 2 methyl phonyl)
	H ₃ C, /=	propionic-acid
	CH ₃ OH	proprome acid
141	Chiral	(R)-3-(4-{3-[4-ethyl-2-
141	(N	(pyridine-2-carbonyl)-
	CH ₃	phenoxy]-butoxy}-2-
	H ₃ C, /=(/=(1	methyl-phenyl)-propionic
		acid
	ČН ₃ ОН	liona .

	Structure	Name
142	⊘ s	3-(2-methyl-4-[3-[2-
		(thiophene-2-carbonyl)-4-
	F = O CH ₃	trifluoromethoxy-phenoxy]
	FO-{_}O-\O	butoxy phenyl) propionic
	CH ₃ OH	acid
143	P _s	3 (4-{3-[4-ethyl-2-
	\\ \\ \\ \\ \\ \\ \\ \\ \\ \\ \\ \\ \\	(thiophene 2 carbonyl)
	O CH ₃	phenoxy] butoxy] 2
	H ₃ C O O	methyl-phenyl) propionic
	сн, он	acid
144	~	3 (4 {3 [4 ethyl-2
		(naphthalene-1-earbonyl)-
		phenoxy]-butoxy}-2-
	H ₃ C	methyl-phenyl)-propionie
	CH ₃ OH	acid
145		3 (4 13 [4 ethyl 2 (1
143		
		phenyl vinyl) phenoxy]
	H ₃ C, /= CH ₃	butoxy 2-methyl-phenyl)
		propionic acid
	cH³OH	
146		3-{4-[3-(2-benzoyl-
		phenoxy) butoxy]-2-
		methyl-phenyl}-propionie
		neid
	CH ₃ OH	
	OH OH	

	Structure	Name
147	H ₃ C—OOOOOOOOOOOOOOOOOOOOOOOOOOOOOOOOOOOO	3-(4-[3-(2-benzoyl-4-methyl-phenoxy) butoxy] 2-methyl-phenyl) propionie aeid
148	H ₃ C CH ₃ O-CH ₃ OH	3-(4-[3-(2-benzyl-4-ethyl- phenoxy)-butoxy]-2- methyl-phenyl}-propionie neid
149	Br—CH ₃ OH	3-(4-[3-(2-benzeyl-4-bromo-phenexy)-butoxy]- 2-methyl-phenyl}-propionie aeid
150	H ₂ C CH ₃ O CH ₃ O O O O O O O O O O O O O O O O O O O	3 (4 [3 (2 benzoyl 4 butyl- phenoxy) butoxy] 2- methyl phenyl propionie aeid
151	H ₃ C CH ₃ OH	3 (4 [3 (2 benzeyl 4- propyl-phenoxy) butoxy] 2 methyl-phenyl] propionie acid
152	CH ₃ OH	3-{4-[4-(2-benzoyl-4-ethyl- phenoxy)-1-methyl- butoxy]-2-methyl-phenyl - propionic-acid

	Structure	Name
1.75	Structure	
153	H ₃ C CH ₃ OH	3-[4-[4-(2-benzoyl-4-ethyl- phenoxy)-pentyloxy]-2- methyl-phenyl}-propionic acid
154	H ₃ C CH ₃ OCH	3-{4-[3-(2-benzoyl-4-ethyl-phenoxyl-2-methyl-propoxyl-2-methyl-phenyl}-propionic-acid
155	H ₃ C CH ₃ OH	3-(4-[3-(2-benzoyl-4-ethyl- phenoxy) propoxy] 2- methyl-phenyl} propionic acid
156	H ₃ C CH ₃	3 (4 3 [4 ethyl-2 (4- fluoro-benzoy!) phenoxy]- propoxy} 2 methyl- phenyl) propionic acid
157	F CH ₃ OO OO OO	3 (4 [3 [4 ethyl 2 (2- trifluoromethyl-benzoyl)- phenoxyl-propoxyl-2- methyl-phenyl) propionie acid

	Structure	Name
158	H ₅ C O O O O O O O O O O O O O O O O O O O	3 (4 [3 [4 ethyl-2 (3-trifluoromethyl-benzey!) phenoxy] propoxy] 2-methyl-phenyl) propionie aeid
159	S CH ₃ COO OO	3 (4 [3 [4 ethyl-2- (thiophene 2 earbonyl)- phenoxy] propoxy] 2- methyl-phenyl) propionie acid
160	H ₃ C CH ₃ O	3 [4 [3 (2 benzyl 4 ethyl- phenoxy) propoxy] 2- methyl-phenyl}-propionie neid
161	H ₃ C OH	3 (4 [3 [4 ethyl-2- (naphthalene 1 carbonyl)- phenoxy] propoxy] 2- methyl-phenyl) propionie acid
162	H ₃ C CH ₃ OH	3 (4 [3 [4 ethyl 2 (1- phenyl-vinyl) phenoxy]- propoxy] 2 methyl- phenyl) propionic acid

	Structure	Name
163	H ₃ C O H ₅ C O O H ₅ C O O O O O O O O O O O O O O O O O O O	2-[4-[3 (2-benzoyl 4-ethyl-phenoxy) butoxy] phenoxyl 2-methyl-propionic-acid
164	H ₃ C	2-[4-[3-(2-benzoy)-4-ethyl-phenoxy)-2-methyl-propionic soid
165	H ₃ C O O O O O O O O O O O O O O O O O O O	2 [4 [3 (2 benzyl 4 ethyl- phenoxy) butoxy]- phenoxyl 2-methyl- propionie aeid
166	Br H ₃ C O O H ₃ C O H ₃ C O H ₃ C O H ₃ C O O O O O O O O O O O O O O O O O O O	2-[4-[3-(2-benzoy]-4-bromo-phenoxy)-butoxy]-phenoxy]-2-methyl-propionic-acid

	A	Lav
	Structure	<u>Name</u>
167	H ₃ C O O O O O O O O O O O O O O O O O O O	2-[4-[3-(2-benzoyl-4-butyl- phenoxy)-butoxy]- phenoxy]-2-methyl- propionic-acid
168	Chral	(R) 3-14 [3 (4-chlore-2- phenoxy-phenoxy)- butoxy] 2-methyl-phenyl}- propionic-acid
169	Chiel Ch	(R) 3 [2 methyl 4 [3 (2 phenoxy 4 trifluoromethyl phenoxy) butoxy] phenyl] propionic acid
170	F O CH ₃	(R) 3-{2-methyl-4-{3-(2-phenoxy-4-trifluoromethoxy-phenoxy-butoxy}-phenoxy-phenoxy-phenoxy}-phenyl}-propionic
171	Chiral O CH ₃ O CH ₃ O O O O O O O O O O O O O O O O O O O	(R) 3 {2-methyl 4 [3 (4-methyl 2-phenoxy-phenoxy) butoxy] phenyl}-propionic-acid

	Structure	Name
172	Chiral	(R)-14-[3-(4-ehloro-2-
		phenoxy-phenoxy)-
	O CH ₃	butoxy] 2 methyl-
	c -{ }-0	phonylsulfanyl)-acetic-acid
	CH ₃ OH	
173	<u></u>	3-{4-{3-(4-chloro-2-
		phenoxy phenoxy)
	OCH ₃	propoxy] 2-methyl-
	c -{\rightarrow-0-{\r	phenyl}-propionie-acid
	ОН	
174	Chiral	(R) 3 (4 [3 (2
	l s	benzo[b]thiophen-3-yl-4-
	сн,	chloro-phenoxy)-butoxy]-2-
	CI	methyl-phenyl}-propionic
	CH ₃ OH	neid
175	Chiral	(R)- 3-{4-[3-(4-chloro-2-
	,cH³	pyridin-3-yl-phenoxy)-
		butoxy]-2-methyl-phenyl}-
	CH ₃ OH	propionic acid
176	Chiral	(R) 3 {4 [3 (4 chloro 2
		phenoxy-phenoxy)-
	Ò	butoxy]-phenyl]-2,2-
	c -{_}-0\0-\	difluoro-propionie acid
	CH ₃ F F OH	
177	Chiral	%)(R) 3 {3 bromo 4 [3 (4-
		chloro 2 phenoxy-
	O Br	phenoxy)-butoxy]-phenyl}-
	c:—(propionie acid
	CH ₃ OH	

	Structure	Name
178	Chval	(R) 3 [4 [3 (4 ehloro-2- phenoxy-phenoxy) butoxy] 3 methyl phenyll- propionic-acid
179	CI—CH ₃ O—CH ₃ HO	(R) {3-bromo 4 {3 (4- ehloro 2 phenoxy- phenoxy) butoxy} pheny}}- neetic acid
180	F-F O CH ₃ Br-CH ₃ O O O O O O O O O O O O O O O O O O O	(R) 3- [4-[3-(4-bromo-2-trifluoromethoxy-phenoxy)-butoxy] 2-methyl-phenyl}-propionic acid
181	CI————————————————————————————————————	(R) [4-[3 (4 chlore 2- phenoxy-phenoxy)- butoxy] 3 methyl-phenyl}- nectic neid
182	Chiral Chiral	(R) [4 [3 (4 chloro 2- phenoxy-phenoxy) butoxy] phenyl acetic acid
183	CI CH ₃ OH OH	(R) 3-[4-[3-(4-ehloro-2- phenoxy-phenoxy)- butoxy]-2-trifluoromethyl- phenyl}-propionie-aeid

	Structure	Name
184	Christ	(R) [4-[3-(4-chloro-2- phenoxy phenoxy) butylsulfanyl] 2-methyl- phenoxy -acetic-acid
185	Chiral Chiral	(R) 3 {4 [3 (4 chloro-2- phenoxy phenoxy) butylsulfanyl]-2-methyl- phenyl}-propionic acid
186	CI OChiral OH	(R) 3-[2-Chloro 4-[3-(4-chloro 2-phenoxy-phenoxy-phenoxy) butoxy] phenyl]-propionic-acid
187	CI Chiral OH	(R) 3 [4 [3 (4 Chloro 2- phenoxy-phenoxy) butoxy] 2 fluoro-phenyl - propionic-acid
188	CI OH	(R) 3-{4-{3-(4-Chloro-2- phenoxy-phenoxy)- butoxy}-2-ethyl-phenyll- propionic-acid
189	CI Chiral OH	(R) 3 {4 [3 (2 Benzoyl 4- ethyl-phenoxy) butoxy] 2- ehloro-phenyl -propionic acid

	Structure	Name
190	F OChiral OH	(R) 3 {4 [3 (2-Benzoyl 4-ethyl-phenoxy) butoxy] 2-fluoro-phenyl -propionie
191	Cl Chiral OH	(R) 3 {4 [3 (4 Chloro 2- phenoxy-phenoxy) butoxy] phenyl}-propionic acid
192	Chiral	(R) 3 {4 [3 (2 Benzoyl 4- ethyl-phenoxy) butoxy]- phenyl}-propionic acid
193	CI Chiral OH	(R) 3 {4 [3 (4 Chloro 2- phenoxy-phenoxy)- pentyloxy] 2-methyl- phenyl} propionic acid
194	OChiral OH	(R) 3 [4 [3 (2 Benzoyl 4- ethyl phenoxy) pentyloxy] 2 methyl phenyl] propionie acid
195	Chiral	(R) [4-[3 (3 Benzeyl- naphthalen 2 yloxy)- butoxy] 2 methyl- phenylsulfanyl] acetic acid

	Structure	Name
196	Chiral	(R) 3-{4-{3-(3-Benzoyl- naphthalen 2-yloxy)- butoxy] 2-methyl-phenyl}- propionic-acid
197	Chiral OH	(R) 3 [4 [3 (4 Ethyl 2- phenoxy-phenoxy)- butylsulfanyl] 2-methyl- phenyl]-propionie acid
198	Chiral	(R) 3 [4 [3 (4 Isopropyl 2- phenoxy phenoxy) butylsulfanyl] 2 methyl phenyl} proptonic acid
199	Cl OH OH	(R) 3 [4 [3 (4 Chloro 2- phenoxy phenoxy) butoxy] 2 propyl phenyl) propionie aeid
200	Cl Chiral	(R) {4 [3 (4-Chlore 2- phenoxy-phenoxy)- butoxy] 2 ethyl- phenylsulfanyl} acetic acid
201	CI Chiral OH	(R) 3-{4-{3-(2-Benzoyl- 4,5-dichloro phenoxy)- butoxy}-2-methyl-phenyl}- propionic-acid

	Structure	Name
202	CF ₃ Chiral OH	(R) 3 {2-Methyl-4-[3-(2- phenoxy-4-trifluoromethyl- phenoxy)-butylsulfanyl]- phenyl}-propionie acid
203	Chiral	(R) 3 {2 Ethyl 4 [3 (4 ethyl 2 phenoxy phenoxy) butoxy] phenyl} propionic acid
204	CF ₅ OH Chiral	(R) 3 {2 Ethyl 4 [3 (2- phenoxy 4 trifluoromethyl- phenoxy)-butoxy]-phenyl}- propionic-acid
205	Chiral	(R) 3 {4 [3 (2 Benzoyl 4 ethyl phenoxy) butoxy] 2-ethyl phenyl} propionie acid
206	CF ₃ OH	(R) 3 [2 Ethyl 4 [1- methyl 3 (2 phenoxy 4- trifluoromethyl phenoxy)- propoxy] phenyl}- propionic acid
207	F F O OH OH	(R) 3 [2-Methyl 4 [1- methyl 3 (2-phenoxy 4- trifluoromethoxy-phenoxy) propylsulfanyl]-phenyl]- propionic-acid

	Structure	Name
208	Chiral OH	(S) 3 {4-{3-(4-Chloro-2- phenoxy-phenoxy)} butoxy} 2 ethyl phenyl} propionic-acid
209	CI OH	3 [4 [3 (4 Chloro 2- phenoxy-phenoxy)- propoxy] 2 ethyl phenyl}- propionic-acid
210	Chiral OH	(R) 3 [4 [3 (2,4 Diphenexy phenexy) butoxy] 2 ethyl phenyll- propionic acid
211	Cl Clis - Isomer 2	2-{4-[4-(4-Chloro-2- phenoxy-phenyl)-3-methyl- butoxy]-2-methyl-phenyl}- cyclopropancearboxylic acid
212	H ₃ C CH ₃ CH ₃ OH	(R, S) 2 [4 [3 (4 Ethyl 2- phenylsulfanyl-phenoxy) butoxy] phenoxy] 2 methyl- propionic acid
213	H ₃ C CH ₃ H ₅ OH CH ₃	2-(4-[3-(R,S-2- Benzenesulfinyl 4-ethyl- phenoxy) butoxy]- 2-methyl-phenylsulfanyl)- 2-methyl-propionic acid (enamtiomer-pair-l)

	Structure	<u>Name</u>
214	F CH ₃ CH ₃ OH	(R, S) 2 [4-[3-(2- Cyclopropylmethyl 4- trifluoromethyl- phenoxy) butoxyl- phenoxyl 2-methyl- propionic acid
215	H ₃ C CH ₃ CH ₃ CH ₃	(R, S) 2 Methyl 2 (4 [3 (2- methyl 3-phenyl 7-propyl- benzofuran-6-yloxy)- butoxy] phenoxy)- propionie acid
216	CH ₃ CH ₅ OH CH ₅	(R, S) 2 Methyl 2 (4 [3 (4- methyl 3 phenyl 7 propyl- benzofuran 6 yloxy)- butoxy] phenoxy]- propionie acid
217	F F CH ₃ H,C OH OH	(R, S) 2 4 3 (2 Cyclopropylmethyl 4 trifluoromethyl- phenoxy) butoxy] 2- methyl-phenoxy] 2-methyl- propionie acid
218	F F CH ₃ OH	(R, S) 3-{4-[3-(2- Cyclopropylmethyl-4- trifluoromethyl- phenoxy) butoxy 2-

	Structure	Name
		methyl-phenyl}-propionie
219	H ₃ C CH ₃ OH	3-[R-4-[3-(R, S-2-) Benzenesulfinyl-4-ethyl- phenoxy)-butoxy]- 2-methyl-phenyl]-propionie acid
220	H ₃ C CH ₃ CH ₃ OH	3 (4 [3 (4 Ethyl 2- phenylsulfanyl phenoxy)- butoxy] 2-methyl-phenyl]-propionie acid-isomer 2
221	H ₃ C CH ₃ CH ₃ OH	(R, S) 2-{4-[3-(4-Ethyl-2- phonylsulfanyl-phonoxy)- butoxy] phonoxy} 2- methyl-propionic acid
222	H ₂ C CH ₃ O OH	(R, S) 3 [44]3 (R, S-2- Benzenesulfinyl 4-ethyl- phenoxy) butoxy] 2-methyl-phenyll- propionic acid
223	H ₃ C CH ₃ CH ₃ OH	(R, S) 2 [4 [3 (R, S 2 Benzenesulfinyl 4-ethyl- phenoxy) -butoxy] 2-methyl- phenoxyl 2-methyl-

	Structure	Name
		propionic acid
224	H ₃ C CH ₃ OH	(R, S) 3 [4 [3 (2- Benzenesulfonyl 4 ethyl- phenoxy)-butoxy]-2- methyl-phenyl]-propionic acid
225	FFF CH ₃ OOH	3- [4-[3-(2-Benzoyl-4- trifluoromethexy-phenoxy)- butoxy] 2-methyl-phenyl}- propionic acid

30. (Withdrawn). The compound of Claim 29, wherein the compound is

or a pharmaceutically acceptable salt, solvate or hydrate thereof.

31. (Previously Presented). A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a compound of Claim 1 or a pharmaceutically acceptable salt, solvate or hydrate thereof.

- 32. (Canceled).
- 33. (Canceled).
- 34. (Canceled).
- 35. (Canceled).
- 36. (Canceled).

- (Canceled).
- 38. (Canceled).
- 39. (Canceled).
- 40. (Canceled).
- 41. (Canceled).
- 42. (Canceled).
- (Previously Presented). A method for lowering blood-glucose in a mammal
 in need thereof comprising the step of administering an effective amount of a compound of Claim
 1.
 - 44. (Canceled).
 - 45. (Canceled).
 - 46. (Canceled).
 - 47. (Canceled).
 - 48. (Canceled).
 - 49. (Canceled)